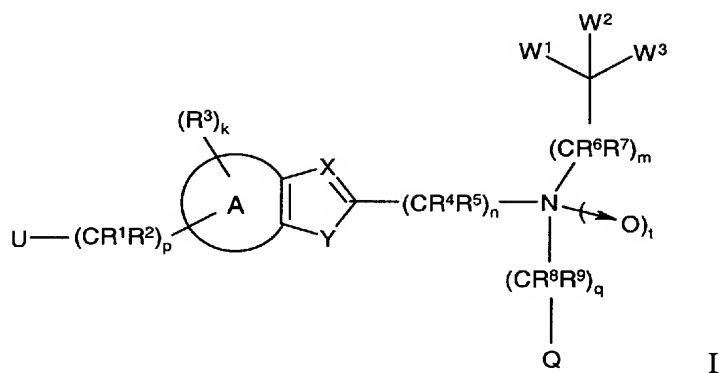


### Amendments to the claims

This listing of claims will replace all prior versions, and listings, of claims in the application:

#### Listing of Claims:

1. (Previously presented): A compound of Formula I:



wherein:

X is CH or N;

Y is N(R<sup>10</sup>), O, or S, wherein t is 0 or 1 when Y is N(R<sup>10</sup>) or O, and t is 0 when Y is S;

U is selected from halo, -OR<sup>10</sup>, -NR<sup>14</sup>R<sup>15</sup>, nitro, cyano, -COOR<sup>10</sup>, -COR<sup>13</sup>, -OCOR<sup>13</sup>, -CONR<sup>14</sup>R<sup>15</sup>, -N(R<sup>14</sup>)COR<sup>13</sup>, -SO<sub>3</sub>H, -SO<sub>2</sub>NR<sup>14</sup>R<sup>15</sup>, -C(=NR<sup>17</sup>)NR<sup>14</sup>R<sup>15</sup>, -N(R<sup>14</sup>)SO<sub>2</sub>R<sup>16</sup>, and a 5 or 6-membered heterocyclic group;

A is a phenyl fused ring moiety or a pyridyl fused ring moiety, wherein when A is a phenyl ring moiety, k is 0-3 and t is 0 or 1 and when A is a pyridyl ring moiety, k is 0-2 and t is 0;

W<sup>1</sup> is selected from C<sub>3</sub>-C<sub>8</sub> cycloalkyl, aryl and Het, wherein said C<sub>3</sub>-C<sub>8</sub> cycloalkyl, Ar and Het are optionally unsubstituted or substituted with one or more groups independently selected from halo, cyano, nitro, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-C(O)SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-COR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>3</sub>H, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OCOR<sup>13</sup>,

-C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)OR<sup>13</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)NR<sup>11</sup>R<sup>12</sup>, and -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, where said  
C<sub>1</sub>-C<sub>6</sub> alkyl, is optionally unsubstituted or substituted by one or more halo  
substituents;

W<sup>2</sup> is selected from H, halo, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>2</sub>-C<sub>6</sub> alkenyl, C<sub>2</sub>-C<sub>6</sub> alkynyl,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-C(O)SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-COR<sup>13</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-OCOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OCONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>CONR<sup>11</sup>R<sup>12</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-Het, -C<sub>0</sub>-C<sub>6</sub> alkyl-Ar and  
-C<sub>0</sub>-C<sub>6</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, wherein said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or  
substituted by one or more halo substituents, and wherein the C<sub>3</sub>-C<sub>7</sub> cycloalkyl, Ar  
and Het moieties of said -C<sub>0</sub>-C<sub>6</sub> alkyl-Het, -C<sub>0</sub>-C<sub>6</sub> alkyl-Ar and  
-C<sub>0</sub>-C<sub>6</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl are optionally unsubstituted or substituted with one or  
more groups independently selected from halo, cyano, nitro, C<sub>1</sub>-C<sub>6</sub> alkyl,  
C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-C(O)SR<sup>10</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-COR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>3</sub>H, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>NR<sup>11</sup>R<sup>12</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OCOR<sup>13</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)OR<sup>13</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)NR<sup>11</sup>R<sup>12</sup>, and -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, where said  
C<sub>1</sub>-C<sub>6</sub> alkyl, is optionally unsubstituted or substituted by one or more halo  
substituents;

W<sup>3</sup> is selected from the group consisting of: H, halo, C<sub>1</sub>-C<sub>6</sub> alkyl,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-C(O)SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-COR<sup>13</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-OCOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OCONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>CONR<sup>11</sup>R<sup>12</sup>,  
-C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-Het, -C<sub>1</sub>-C<sub>6</sub> alkyl-Ar and  
-C<sub>1</sub>-C<sub>6</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, wherein said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or  
substituted by one or more halo substituents;

Q is selected from C<sub>3</sub>-C<sub>8</sub> cycloalkyl, Ar and Het; wherein said  
C<sub>3</sub>-C<sub>8</sub> cycloalkyl, Ar and Het are optionally unsubstituted or substituted with one or

more groups independently selected from halo, cyano, nitro, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-C(O)SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-COR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>3</sub>H, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OCOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)NR<sup>11</sup>R<sup>12</sup>, and -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, where said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents;

p is 0-8;

n is 2;

m is 0 or 1;

q is 0 or 1;

t is 0 or 1;

each R<sup>1</sup> and R<sup>2</sup> are independently selected from H, halo, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SR<sup>10</sup>, -C<sub>1</sub>-C<sub>6</sub> alkyl-Het, -C<sub>1</sub>-C<sub>6</sub> alkyl-Ar and -C<sub>1</sub>-C<sub>6</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, or R<sup>1</sup> and R<sup>2</sup> together with the carbon to which they are attached form a 3-5 membered carbocyclic or heterocyclic ring, wherein said heterocyclic ring contains one, or more heteroatoms selected from N, O, and S, where said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents;

each R<sup>3</sup> is the same or different and is independently selected from halo, cyano, nitro, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-Ar, -C<sub>0</sub>-C<sub>6</sub> alkyl-Het, -C<sub>0</sub>-C<sub>6</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-C(O)SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-COR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>3</sub>H, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-SOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OCOR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-OC(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>C(O)NR<sup>11</sup>R<sup>12</sup>, and -C<sub>0</sub>-C<sub>6</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, wherein said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents;

each  $R^4$  and  $R^5$  is independently selected from H, halo,  $C_1$ - $C_6$  alkyl,  $-C_0$ - $C_6$  alkyl-Het,  $-C_0$ - $C_6$  alkyl-Ar and  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl;

$R^6$  and  $R^7$  are each independently selected from H, halo,  $C_1$ - $C_6$  alkyl,  $-C_0$ - $C_6$  alkyl-Het,  $-C_0$ - $C_6$  alkyl-Ar and  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl;

$R^8$  and  $R^9$  are each independently selected from H, halo,  $C_1$ - $C_6$  alkyl,  $-C_0$ - $C_6$  alkyl-Het,  $-C_0$ - $C_6$  alkyl-Ar and  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl;

$R^{10}$  is selected from H,  $C_1$ - $C_6$  alkyl,  $C_3$ - $C_6$  alkenyl,  $C_3$ - $C_6$  alkynyl,  $-C_0$ - $C_6$  alkyl-Ar,  $-C_0$ - $C_6$  alkyl-Het and  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl;

each  $R^{11}$  and each  $R^{12}$  are independently selected from H,  $C_1$ - $C_6$  alkyl,  $C_3$ - $C_6$  alkenyl,  $C_3$ - $C_6$  alkynyl,  $-C_0$ - $C_6$  alkyl-Ar,  $-C_0$ - $C_6$  alkyl-Het and  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl, or  $R^{11}$  and  $R^{12}$  together with the nitrogen to which they are attached form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S;

$R^{13}$  is selected from  $C_1$ - $C_6$  alkyl,  $C_3$ - $C_6$  alkenyl,  $C_3$ - $C_6$  alkynyl,  $-C_0$ - $C_6$  alkyl-Ar,  $-C_0$ - $C_6$  alkyl-Het and  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl;

$R^{14}$  and  $R^{15}$  are each independently selected from H,  $C_1$ - $C_6$  alkyl,  $C_3$ - $C_6$  alkenyl,  $C_3$ - $C_6$  alkynyl,  $-C_0$ - $C_6$  alkyl-Ar,  $-C_0$ - $C_6$  alkyl-Het,  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl,  $-C_0$ - $C_6$  alkyl-O-Ar,  $-C_0$ - $C_6$  alkyl-O-Het,  $-C_0$ - $C_6$  alkyl-O- $C_3$ - $C_7$  cycloalkyl,  $-C_0$ - $C_6$  alkyl-S(O)<sub>x</sub>- $C_1$ - $C_6$  alkyl,  $-C_0$ - $C_6$  alkyl-S(O)<sub>x</sub>-Ar,  $-C_0$ - $C_6$  alkyl-S(O)<sub>x</sub>-Het,  $-C_0$ - $C_6$  alkyl-S(O)<sub>x</sub>- $C_3$ - $C_7$  cycloalkyl,  $-C_0$ - $C_6$  alkyl-NH-Ar,  $-C_0$ - $C_6$  alkyl-NH-Het,  $-C_0$ - $C_6$  alkyl-NH- $C_3$ - $C_7$  cycloalkyl,  $-C_0$ - $C_6$  alkyl-N( $C_1$ - $C_4$  alkyl)-Ar,  $-C_0$ - $C_6$  alkyl-N( $C_1$ - $C_4$  alkyl)-Het,  $-C_0$ - $C_6$  alkyl-N( $C_1$ - $C_4$  alkyl)- $C_3$ - $C_7$  cycloalkyl,  $-C_0$ - $C_6$  alkyl-Ar,  $-C_0$ - $C_6$  alkyl-Het and  $-C_0$ - $C_6$  alkyl- $C_3$ - $C_7$  cycloalkyl, where x is 0, 1 or 2, or  $R^{14}$  and  $R^{15}$ , together with the nitrogen to which they are attached, form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S, wherein said  $C_1$ - $C_6$  alkyl is optionally substituted by one or more of the substituents independently selected from the group halo, -OH, -SH, -NH<sub>2</sub>, -NH(unsubstituted  $C_1$ - $C_6$  alkyl), -N(unsubstituted  $C_1$ - $C_6$  alkyl)(unsubstituted  $C_1$ - $C_6$  alkyl), unsubstituted -OC<sub>1</sub>- $C_6$  alkyl, -CO<sub>2</sub>H, -CO<sub>2</sub>(unsubstituted  $C_1$ - $C_6$  alkyl), -CONH<sub>2</sub>, -CONH(unsubstituted  $C_1$ - $C_6$  alkyl), -CON(unsubstituted  $C_1$ - $C_6$  alkyl)(unsubstituted

C<sub>1</sub>-C<sub>6</sub> alkyl), -SO<sub>3</sub>H, -SO<sub>2</sub>NH<sub>2</sub>, -SO<sub>2</sub>NH(unsubstituted C<sub>1</sub>-C<sub>6</sub> alkyl) and -SO<sub>2</sub>N(unsubstituted C<sub>1</sub>-C<sub>6</sub> alkyl)(unsubstituted C<sub>1</sub>-C<sub>6</sub> alkyl);

R<sup>16</sup> is C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-Ar or -C<sub>0</sub>-C<sub>6</sub> alkyl-Het; and

R<sup>17</sup> is H, C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>6</sub> alkyl-Ar or -C<sub>0</sub>-C<sub>6</sub> alkyl-Het;

or a pharmaceutically acceptable salt or solvate thereof.

2. (Original): The compound according to claim 1, wherein p is 0, 1 or 2.

3. (Original): The compound according to claim 1, wherein t is 0.

4. (Previously presented): The compound according to claim 1, wherein R<sup>1</sup> and R<sup>2</sup> are each H.

5. (Previously presented): The compound according to claim 1, wherein A is a phenyl fused ring.

6. (Previously presented): The compound according to claim 1, wherein k is 0.

7. (Previously presented): The compound according to claim 1, wherein U is -OR<sup>10</sup>, -COOR<sup>10</sup>, -CONR<sup>11</sup>R<sup>12</sup> or -NR<sup>11</sup>R<sup>12</sup>.

8. (Previously presented): The compound according to claim 1, wherein U is -OH, -COOH, -CONH<sub>2</sub>, -CON(H)CH<sub>2</sub>-furan-2-yl, or -N(H)CH<sub>2</sub>-furan-2-yl.

Claims 9-10 (Cancelled).

11. (Previously presented): The compound according to claim 1, wherein q is 1.

12. (Previously presented): The compound according to claim 1, wherein  $R^8$  and  $R^9$  are each H.

13. (Previously presented): The compound according to claim 1, wherein Q is a substituted phenyl group, containing one or two substituents selected from halo,  $C_1$ - $C_4$  alkoxy; and  $C_1$ - $C_4$  alkyl or Q is a 1,3-benzodioxolyl or dihydrobenzofuranyl group.

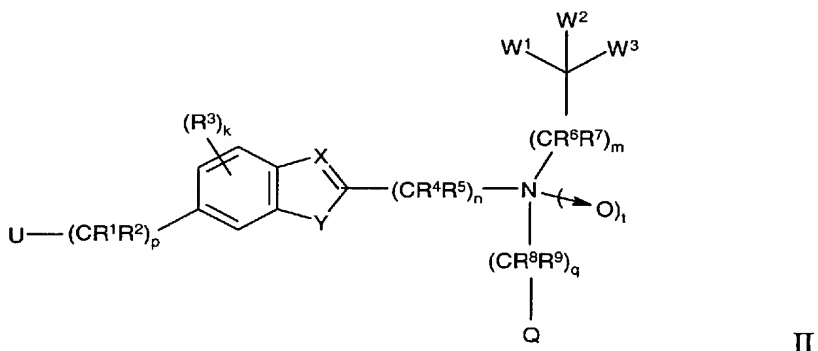
14. (Previously presented): The compound according to claim 1, wherein Q is a phenyl group substituted by one or two substituents selected from chloro, trifluoromethyl and methoxy or is a 1,3-benzodioxolyl or a dihydrobenzofuranyl group.

15. (Previously presented): The compound according to claim 1, wherein m is 1 and  $R^6$  and  $R^7$  are both H.

16. (Previously presented): The compound according to claim 1, wherein  $W^3$  is H

17. (Previously presented): The compound according to claim 1 wherein  $W^1$  and  $W^2$  are each unsubstituted phenyl or  $W^1$  is unsubstituted phenyl and  $W^2$  is methyl.

18. (Previously presented): A compound of Formula II:



wherein:

X is CH or N;

Y is O, or S;

U is selected from halo,  $-OR^{10}$ ,  $-NR^{14}R^{15}$ , cyano,  $-COOR^{10}$ ,  $-OCOR^{13}$ ,  $-CONR^{14}R^{15}$ ,  $-N(R^{14})COR^{13}$ ,  $-SO_2NR^{14}R^{15}$ ,  $-C(=NH)NR^{14}R^{15}$ , and a 5 or 6-membered heterocyclic group;

$W^1$  is selected from  $C_3$ - $C_8$  cycloalkyl, aryl and Het, wherein said  $C_3$ - $C_8$  cycloalkyl, Ar and Het are optionally unsubstituted or substituted with one or more groups independently selected from halo, cyano, nitro,  $C_1$ - $C_6$  alkyl,  $C_3$ - $C_6$  alkenyl,  $C_3$ - $C_6$  alkynyl,  $-C_0$ - $C_4$  alkyl- $CO_2R^{10}$ ,  $-C_0$ - $C_4$  alkyl- $C(O)SR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $CONR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $COR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $NR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $SR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $OR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $SO_3H$ ,  $-C_0$ - $C_4$  alkyl- $SO_2NR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $SO_2R^{10}$ ,  $-C_0$ - $C_4$  alkyl- $SOR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $OCOR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $OC(O)NR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $OC(O)OR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $NR^{11}C(O)OR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $NR^{11}C(O)NR^{11}R^{12}$ , and  $-C_0$ - $C_4$  alkyl- $NR^{11}COR^{13}$ , where said  $C_1$ - $C_6$  alkyl is optionally unsubstituted or substituted by one or more halo substituents;

$W^2$  is selected from H, halo,  $C_1$ - $C_6$  alkyl,  $C_2$ - $C_6$  alkenyl,  $C_2$ - $C_6$  alkynyl,  $-C_0$ - $C_4$  alkyl- $NR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $SR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $OR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $CO_2R^{10}$ ,  $-C_0$ - $C_4$  alkyl- $C(O)SR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $CONR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $COR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $OCOR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $OCONR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $NR^{11}CONR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $NR^{11}COR^{13}$ ,  $-C_0$ - $C_4$  alkyl-Het,  $-C_0$ - $C_4$  alkyl-Ar and  $-C_0$ - $C_4$  alkyl- $C_3$ - $C_7$  cycloalkyl, wherein said  $C_1$ - $C_6$  alkyl is optionally unsubstituted or substituted by one or more halo substituents, and wherein the  $C_3$ - $C_7$  cycloalkyl, Ar and Het moieties of said  $-C_0$ - $C_4$  alkyl-Het,  $-C_0$ - $C_4$  alkyl-Ar and  $-C_0$ - $C_4$  alkyl- $C_3$ - $C_7$  cycloalkyl are optionally unsubstituted or substituted with one or more groups independently selected from halo, cyano, nitro,  $C_1$ - $C_6$  alkyl,  $C_3$ - $C_6$  alkenyl,  $C_3$ - $C_6$  alkynyl,  $-C_0$ - $C_4$  alkyl- $CO_2R^{10}$ ,  $-C_0$ - $C_4$  alkyl- $C(O)SR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $CONR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $COR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $NR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $SR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $OR^{10}$ ,  $-C_0$ - $C_4$  alkyl- $SO_3H$ ,  $-C_0$ - $C_4$  alkyl- $SO_2NR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $SO_2R^{10}$ ,  $-C_0$ - $C_4$  alkyl- $SOR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $OCOR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $OC(O)NR^{11}R^{12}$ ,  $-C_0$ - $C_4$  alkyl- $OC(O)OR^{13}$ ,  $-C_0$ - $C_4$  alkyl- $NR^{11}C(O)OR^{13}$ ,

-C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>C(O)NR<sup>11</sup>R<sup>12</sup>, and -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, where said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents;

W<sup>3</sup> is selected from the group consisting of: H, halo, C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-C(O)SR<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-COR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OCOR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OCONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-Het, -C<sub>1</sub>-C<sub>4</sub> alkyl-Ar and -C<sub>1</sub>-C<sub>4</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, wherein said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents;

Q is Ar or Het; wherein said Ar and Het are optionally unsubstituted or substituted with one or more groups independently selected from halo, cyano, nitro, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-CO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-C(O)SR<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-CONR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-COR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-SR<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-SO<sub>3</sub>H, -C<sub>0</sub>-C<sub>4</sub> alkyl-SO<sub>2</sub>NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-SO<sub>2</sub>R<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-SOR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OCOR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OC(O)NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OC(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>C(O)OR<sup>13</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>C(O)NR<sup>11</sup>R<sup>12</sup>, and -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>COR<sup>13</sup>, where said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents,

p is 0-4;

n is 2;

m is 0 or 1;

q is 0 or 1;

t is 0;

each R<sup>1</sup> and R<sup>2</sup> are independently selected from H, fluoro, C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-OR<sup>10</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-SR<sup>10</sup>, -C<sub>1</sub>-C<sub>4</sub> alkyl-Het, -C<sub>1</sub>-C<sub>4</sub> alkyl-Ar and -C<sub>1</sub>-C<sub>4</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, where said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents;

each R<sup>3</sup> is the same or different and is independently selected from halo, cyano, C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-NR<sup>11</sup>R<sup>12</sup>, -C<sub>0</sub>-C<sub>4</sub> alkyl-OR<sup>10</sup>,



-C<sub>0</sub>-C<sub>4</sub> alkyl-SO<sub>2</sub>NR<sup>11</sup>R<sup>12</sup>, and -C<sub>0</sub>-C<sub>4</sub> alkyl-CO<sub>2</sub>H, wherein said C<sub>1</sub>-C<sub>6</sub> alkyl is optionally unsubstituted or substituted by one or more halo substituents;

each R<sup>4</sup> and R<sup>5</sup> is independently selected from H, fluoro and C<sub>1</sub>-C<sub>6</sub> alkyl;

R<sup>6</sup> and R<sup>7</sup> are each independently selected from H, fluoro and C<sub>1</sub>-C<sub>6</sub> alkyl;

R<sup>8</sup> and R<sup>9</sup> are each independently selected from H, fluoro and C<sub>1</sub>-C<sub>6</sub> alkyl;

R<sup>10</sup> is selected from H, C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-Het and -C<sub>0</sub>-C<sub>4</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl;

each R<sup>11</sup> and each R<sup>12</sup> are independently selected from H, C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-Het and -C<sub>0</sub>-C<sub>4</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, or R<sup>11</sup> and R<sup>12</sup> together with the nitrogen to which they are attached form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S;

R<sup>13</sup> is selected from C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-Het and -C<sub>0</sub>-C<sub>4</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl;

R<sup>14</sup> and R<sup>15</sup> are each independently selected from H, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-Het, -C<sub>0</sub>-C<sub>4</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-O-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-O-Het, -C<sub>0</sub>-C<sub>4</sub> alkyl-O-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-S(O)<sub>x</sub>-C<sub>1</sub>-C<sub>6</sub> alkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-S(O)<sub>x</sub>-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-S(O)<sub>x</sub>-Het, -C<sub>0</sub>-C<sub>4</sub> alkyl-S(O)<sub>x</sub>-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-NH-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-NH-Het, -C<sub>0</sub>-C<sub>4</sub> alkyl-NH-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-N(C<sub>1</sub>-C<sub>4</sub> alkyl)-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-N(C<sub>1</sub>-C<sub>4</sub> alkyl)-Het, -C<sub>0</sub>-C<sub>4</sub> alkyl-N(C<sub>1</sub>-C<sub>4</sub> alkyl)-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, -C<sub>0</sub>-C<sub>4</sub> alkyl-Ar, -C<sub>0</sub>-C<sub>4</sub> alkyl-Het and -C<sub>0</sub>-C<sub>4</sub> alkyl-C<sub>3</sub>-C<sub>7</sub> cycloalkyl, where x is 0, 1 or 2, or R<sup>14</sup> and R<sup>15</sup>, together with the nitrogen to which they are attached, form a 4-7 membered heterocyclic ring which optionally contains one or more additional heteroatoms selected from N, O, and S, wherein said C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl are optionally substituted by one or more of the substituents independently selected from the group halo, -OH, -SH, -NH<sub>2</sub>, -NH(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl), -N(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl)(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl), unsubstituted -OC<sub>1</sub>-C<sub>4</sub> alkyl, -CO<sub>2</sub>H, -CO<sub>2</sub>(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl), -CONH<sub>2</sub>, -CONH(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl), -CON(unsubstituted

C<sub>1</sub>-C<sub>4</sub> alkyl)(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl), -SO<sub>3</sub>H, -SO<sub>2</sub>NH<sub>2</sub>, -SO<sub>2</sub>NH(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl) and -SO<sub>2</sub>N(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl)(unsubstituted C<sub>1</sub>-C<sub>4</sub> alkyl);  
or a pharmaceutically acceptable salt or solvate thereof.

19. (Previously presented): The compound according to claim 1, wherein: R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>8</sup> and R<sup>9</sup> are each H; U is -OR<sup>10</sup>, -COOR<sup>10</sup>, -CONR<sup>11</sup>R<sup>12</sup> or -NR<sup>11</sup>R<sup>12</sup>; A is a phenyl fused ring; Q is a substituted phenyl group containing one or two substituents selected from halo, C<sub>1</sub>-C<sub>4</sub> alkoxy and C<sub>1</sub>-C<sub>4</sub> alkyl or Q is a 1,3-benzodioxolyl or a dihydrobenzofuranyl group; p is 1 or 2; n is 2; m is 1; q is 1; k is 0; t is 0; W<sup>1</sup> is aryl; W<sup>2</sup> is aryl or C<sub>1</sub>-C<sub>4</sub> alkyl; and W<sup>3</sup> is H; or a pharmaceutically acceptable salt or solvate thereof.

20. (Previously presented): The compound according to claim 1, wherein R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>6</sup>, R<sup>7</sup>, R<sup>8</sup>, R<sup>9</sup> and W<sup>3</sup> are each H; U is -OH, -COOH, -CONH<sub>2</sub>, -CON(H)CH<sub>2</sub>-furan-2-yl, -N(H)CH<sub>2</sub>-furan-2-yl; A is a phenyl fused ring; Q is a phenyl group substituted by one or two substituents selected from chloro, trifluoromethyl and methoxy or Q is a 1,3-benzodioxolyl or a dihydrobenzofuranyl group; p is 1 or 2; n is 2; m is 1; q is 1; k is 0; t is 0; W<sup>1</sup> is unsubstituted phenyl; and W<sup>2</sup> is methyl or unsubstituted phenyl; or a pharmaceutically acceptable salt or solvate thereof.

21. (Currently amended): A compound selected from:  
2-[2-{ [2-chloro-3-(trifluoromethyl)-benzyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid,  
2-[2-{ [2,4-dimethoxy-benzyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid,  
2-[2-{ [(2,3-methylenedioxy)benzyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid,  
2-[2-{ [(2,3-dihydrobenzo[*b*]furan)methyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid,

2-[2-{[4-methoxy-benzyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran  
acetic acid,

(*R*)-2-[2-{[2-chloro-3-(trifluoromethyl)-benzyl] (2-methyl-2-  
phenylethyl)amino }ethyl]-5-benzofuran acetic acid,

(*R*)-2-[2-{[(2,3-dihydrobenzo[*b*]furan)methyl](2-methyl-2-  
phenylethyl)amino }ethyl]-5-benzofuran acetic acid,

(*S*)-2-[2-{[2-chloro-3-(trifluoromethyl)-benzyl] (2-methyl-2-  
phenylethyl)amino }ethyl]-benzofuran acetic acid,

(*S*)-2-[2-{[(2,3-dihydrobenzo[*b*]furan)methyl](2-methyl-2-  
phenylethyl)amino }ethyl]-5-benzofuran acetic acid,

2-{2-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]-ethyl}-6-  
benzofuran acetic acid,

2-[2-{ [(2,3-methylenedioxy)benzyl](2,2-diphenylethyl)amino }ethyl]-6-  
benzofuran acetic acid,

2-[2-{ [(2,4-dimethoxy)benzyl](2,2-diphenylethyl)amino }ethyl]-6-benzofuran  
acetic acid,

2-{2-[(4-methoxy-benzyl)(2,2-diphenylethyl)amino]-ethyl}-6-benzofuran  
acetic acid,

2-{2-[(2-chloro-3-(trifluoromethyl)-benzyl)-(2,2-diphenylethyl-amino] ethyl}-  
benzofuran-6-yl)-N-furan-2-yl methyl-acetamide,

2-{2-[(2,4-dimethoxy-benzyl)(2,2-diphenylethyl)-amino]ethyl}-benzofuran-6-  
yl)-N-furan-2-yl methyl -acetamide,

2-{2-[(2-chloro-3-(trifluoromethyl)-benzyl) (2,2-diphenylethyl-amino]ethyl}-  
benzofuran-6-yl)-acetamide,

2-{3-[(2-chloro-3-( trifluoromethyl)-benzyl)-(2-phenyl-propyl)-amino]-  
propyl}-benzofuran-6-yl)-acetic acid,

2-(2-{3-[(2-chloro-3-(trifluoromethyl)-benzyl-(2,2-diphenylethyl)-amino]-  
propyl}-benzofuran-6-yl)-ethanol,

2-(2-{3-[(2,4-dimethoxy)-benzyl-(2,2-diphenylethyl)-amino]-propyl}-  
benzofuran-6-yl)-ethanol,

2-{3-[(2-chloro-3-(trifluoromethyl)-benzyl)-((*R*)-2-phenyl-propyl)-amino]-propyl}-benzofuran-6-yl)-acetic acid,

2-{3-[(2-chloro-3-(trifluoromethyl)-benzyl)-((*S*)-2-phenyl-propyl)-amino]-propyl}-benzofuran-6-yl)-acetic acid,

(2-chloro-3-trifluoromethyl-benzyl)-(2,2-diphenyl-ethyl)-[3-(6-{2-[(furan-2-ylmethyl)-amino]-ethyl-benzofuran-2-yl)-propyl]-amine,

~~and a stereoisomer, a stereoisomeric mixture or racemate thereof~~ and a pharmaceutically acceptable salt or solvate thereof.

22. (Currently amended): The compound according to claim 21, selected from:

2-[2-{ [2,4-dimethoxy-benzyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid,

(*R*)-2-[2-{ [(2,3-dihydrobenzo[*b*]furan)methyl](2-methyl-2-phenylethyl)amino }ethyl]-5-benzofuran acetic acid,

2-{2-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]-ethyl}-6-benzofuran acetic acid,

2-[2-{ [(2,4-dimethoxy)benzyl](2,2-diphenylethyl)amino }ethyl]-6-benzofuran acetic acid,

~~and a stereoisomer, a stereoisomeric mixture or racemate thereof~~ and a pharmaceutically acceptable salt or solvate thereof.

23. (Previously presented): A pharmaceutical composition comprising a compound according to claim 1 and a pharmaceutically acceptable carrier or diluent.

Claims 24-26. (Cancelled).

27. (Previously presented) A method for the prevention or treatment of an LXR mediated disease or condition comprising administering a therapeutically

effective amount of a compound according to claim 1, wherein said LXR mediated disease or condition is atherosclerosis.

28. (Previously presented): A method for the prevention or treatment of an LXR mediated disease or condition comprising administering a therapeutically effective amount of a compound according to claim 1, wherein said LXR mediated disease or condition is inflammation.

29. (Previously presented): A method for increasing reverse cholesterol transport, said method comprising administering a therapeutically effective amount of a compound according to claim 1.

30. (Previously presented): A method for inhibiting cholesterol absorption, said method comprising administering a therapeutically effective amount of a compound according to claim 1.

Claims 31-38 (Cancelled).

39. (Currently amended): A compound selected from the group:

2-[ 2-[( 2,2-diphenylethyl)amino]ethyl]-5-benzofuran acetic acid methyl ester,  
2-[2-[[2-chloro-3-(trifluoromethyl)benzyl-(2,2-diphenylethyl)amino]ethyl]-5-benzofuran acetic acid methyl ester,

2-[2-{ [2,4-dimethoxy-benzyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid methyl ester,

2-[2- { [(2,3-methylenedioxy)benzyl](2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid methyl ester,

2-[2-{ [(2,3-dihydrobenzo[*b*]furan)methyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid methyl ester,

2-[2-{ [4-methoxy-benzyl] (2,2-diphenylethyl)amino }ethyl]-5-benzofuran acetic acid methyl ester,

(*R*)-2-[ 2-[(2-methyl-2-phenylethyl)amino]ethyl]-5-benzofuran acetic acid methyl ester,

(*R*)-2-[2-{ [2-chloro-3-(trifluoromethyl)-benzyl] (2-methyl-2-phenylethyl)amino }ethyl]-5-benzofuran acetic acid methyl ester,

(*R*)-2-[2-{ [(2,3-dihydrobenzo[*b*]furan)methyl](2-methyl-2-phenylethyl)amino }ethyl]-5-benzofuran acetic acid methyl ester,

(*S*)-2-[ 2-[(2-methyl-2-phenylethyl)amino]ethyl]-5-benzofuran acetic acid methyl ester,

(*S*)-2-[2-{ [2-chloro-3-(trifluoromethyl)-benzyl] (2-methyl-2-phenylethyl)amino }ethyl]-benzofuran acetic acid methyl ester,

(*S*)-2-[2-{ [(2,3-dihydrobenzo[*b*]furan)methyl](2-methyl-2-phenylethyl)amino }ethyl]-5-benzofuran acetic acid methyl ester,

2-{ 2-[(2,2-diphenylethyl)amino]-ethyl }-6-benzofuran acetic acid methyl ester,

2-{ 2-[[2-chloro-3-(trifluoromethyl)benzyl](2,2-diphenylethyl)amino]-ethyl }-6-benzofuran acetic acid methyl ester,

2-[2-{ [(2,3-methylenedioxy)benzyl](2,2-diphenylethyl)amino }ethyl]-6-benzofuran acetic acid methyl ester,

2-[2-{ [(2,4-dimethoxy)benzyl](2,2-diphenylethyl)amino }ethyl]-6-benzofuran acetic acid methyl ester,

2-{ 2-[(4-methoxy-benzyl)(2,2-diphenylethyl)amino]-ethyl }-6-benzofuran acetic acid methyl ester,

~~and a stereoisomer, a stereoisomeric mixture or racemate thereof~~ and a pharmaceutically acceptable salt or solvate thereof.